

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM122018\
 Data File : BM018338.D
 Acq On : 20 Dec 2018 19:57
 Operator : JU/SJ
 Sample : J6389-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 P001-SS012-0612-01

Quant Time: Dec 21 01:26:31 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM121518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Dec 15 21:39:37 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.48	152	158175	20.00	ng	-0.02
21) Naphthalene-d8	10.25	136	561977	20.00	ng	-0.02
39) Acenaphthene-d10	14.14	164	267722	20.00	ng	-0.02
64) Phenanthrene-d10	16.91	188	550139	20.00	ng	-0.01
76) Chrysene-d12	21.14	240	691843	20.00	ng	0.00
87) Perylene-d12	23.32	264	729513	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.13	112	1073931	113.42	ng	0.00
7) Phenol-d6	6.69	99	1369444	104.05	ng	-0.01
23) Nitrobenzene-d5	8.64	82	881280	80.44	ng	-0.02
42) 2,4,6-Tribromophenol	15.66	330	391838	99.72	ng	-0.01
45) 2-Fluorobiphenyl	12.75	172	1529116	73.98	ng	-0.02
79) Terphenyl-d14	19.57	244	1976129	64.60	ng	0.00
Target Compounds						
10) Phenol	6.72	94	102555	7.358	ng	94
50) Dimethylphthalate	13.61	163	183002	8.222	ng	# 96
71) Phenanthrene	16.95	178	69231	2.317	ng	# 89
78) Pyrene	19.35	202	91958	2.231	ng	# 86
92) Benzo(g,h,i)perylene	26.12	276	91118	2.670	ng	# 82

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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